

Tocopherol search.

Trying 31060000009999...Open

DIALOG INFORMATION SERVICES

PLEASE LOGON:

***** HHHHHHHH SSSSSSSS? ### Status: Signing onto Dialog *****

ENTER PASSWORD:

***** HHHHHHHH SSSSSSSS? *****

Welcome to DIALOG

Status: Login successfulDialog level 05.12.03D

Last logoff: 02oct06 10:32:10

Logon file001 31oct06 07:53:11

*** ANNOUNCEMENTS ***

NEW FILES RELEASED

***Verdict Market Research (File 769)

***EMCare (File 45)

***Trademarkscan - South Korea (File 655)

***Regulatory Affairs Journals (File 183)

***Index Chemicus (File 302)

***Inspec (File 202)

RESUMED UPDATING

***File 141, Reader's Guide Abstracts

RELOADS COMPLETED

***File 11, PsycInfo

***File 531, American Business Directory

*** The 2005 reload of the CLAIMS files (Files 340, 341, 942)

is now available online.

DATABASES REMOVED

***File 196, FINDEX

***File 468, Public Opinion Online (POLL)

Chemical Structure Searching now available in Prous Science Drug Data Report (F452), Prous Science Drugs of the Future (F453), IMS R&D Focus (F445/955), Pharmaprojects (F128/928), Beilstein Facts (F390), Derwent Chemistry Resource (F355) and Index Chemicus (File 302).

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>>><http://www.dialog.com/whatsnew/>. You can find news about<<<

>>>a specific database by entering HELP NEWS <file number>.<<<

* * *

File 1:ERIC 1966-2006/Sep
(c) format only 2006 Dialog

Set Items Description

--- -----

Cost is in DialUnits

?

Terminal set to DLINK

? b 73,34,399,155,5

31oct06 07:54:43 User291725 Session D6.1

\$0.41 0.116 DialUnits File1

\$0.41 Estimated cost File1

\$0.53 TELNET
\$0.94 Estimated cost this search
\$0.94 Estimated total session cost 0.116 DialUnits

SYSTEM:OS - DIALOG OneSearch

File 73:EMBASE 1974-2006/Oct 27

(c) 2006 Elsevier B.V.

File 34:SciSearch(R) Cited Ref Sci 1990-2006/Oct W4

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File 399:CA SEARCH(R) 1967-2006/UD=14519

(c) 2006 American Chemical Society

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IPCR/8 classification codes now searchable as IC=. See HELP NEWSIPCR.

File 155:MEDLINE(R) 1950-2006/Oct 27

(c) format only 2006 Dialog

File 5:Biosis Previews(R) 1969-2006/Oct W4

(c) 2006 The Thomson Corporation

Set Items Description

? s (alkaloid or morphine or atropine or opioid) same (electron(w)transfer(w)
agent)

>>>Invalid syntax

?

PLEASE ENTER A COMMAND OR BE LOGGED OFF IN 5 MINUTES

? b 73, 34, 399, 155, 5

31oct06 08:00:47 User291725 Session D6.2

\$0.48 0.043 DialUnits File73

\$0.48 Estimated cost File73

\$1.01 0.043 DialUnits File34

\$1.01 Estimated cost File34

\$0.54 0.043 DialUnits File399

\$0.54 Estimated cost File399

\$0.15 0.043 DialUnits File155

\$0.15 Estimated cost File155

\$0.26 0.043 DialUnits File5

\$0.26 Estimated cost File5

OneSearch, 5 files, 0.214 DialUnits FileOS

\$1.86 TELNET

\$4.30 Estimated cost this search

\$5.24 Estimated total session cost 0.330 DialUnits

SYSTEM:OS - DIALOG OneSearch

File 73:EMBASE 1974-2006/Oct 27

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File 34:SciSearch(R) Cited Ref Sci 1990-2006/Oct W4

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File 399:CA SEARCH(R) 1967-2006/UD=14519

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IPCR/8 classification codes now searchable as IC=. See HELP NEWSIPCR.

File 155:MEDLINE(R) 1950-2006/Oct 27

(c) format only 2006 Dialog

File 5:Biosis Previews(R) 1969-2006/Oct W4

(c) 2006 The Thomson Corporation

Set Items Description

? s morphine or atropine or opioid or quinine (w)electron (w) transfer (w)
agent

198329 MORPHINE
136139 ATROPINE
180371 OPIOID
29908 QUININE
3365405 ELECTRON
1480652 TRANSFER
2112835 AGENT
0 QUININE (W) ELECTRON (W) TRANSFER (W) AGENT
S1 457763 MORPHINE OR ATROPINE OR OPIOID OR QUININE (W) ELECTRON (W)
TRANSFER (W) AGENT

? tocopherol(w)phosphate

>>>'OCOPHEROL' not recognized as set or accession number

? tocol(w)phosphate

>>>'OCOL' not recognized as set or accession number

? s quinol or tocopherol(w)phosphate

7837 QUINOL
91705 TOCOPHEROL
988809 PHOSPHATE
141 TOCOPHEROL (W) PHOSPHATE
S2 7978 QUINOL OR TOCOPHEROL (W) PHOSPHATE

? S s1 and s2

457763 S1
7978 S2

S3 17 S1 AND S2

? rd

S4 12 RD (unique items)

? sort/all

>>>Missing range and/or sort tags, please try again.

? sort s4/all/py

S5 12 Sort S4/ALL/PY

? t s5/full/all

5/9/1 (Item 1 from file: 73)

DIALOG(R) File 73:EMBASE

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00428903 EMBASE No: 1975201311

**The prevention by sulphydryl compounds of the toxicity in the cat of 2,6
dimethoxyphenol and its morpholinopropionyl ester**

Loveless A.H.; Maxwell D.R.

Pharmacol. Res. Lab., May Baker Ltd, Dagenham United Kingdom

BRIT.J.PHARMACOL. 1975, 53/1 (93-98)

CODEN: BJPCA

DOCUMENT TYPE: Journal

LANGUAGE: ENGLISH

Intravenous (-) 2,6 dimethoxyphenyl 2 morpholinopropionate hydrochloride (M&B 16.573) produced anaesthesia of short duration in the mouse, rat, rabbit, cat, dog and monkey. In the cat but not in other species, a severe and usually fatal toxic reactions was seen 1-2 hr after administration. This toxic reaction but not the anaesthetic properties of M&B 16.573 was prevented by the intravenous administration of cysteine or N acetylcysteine. Cysteamine or dimercaprol were ineffective. Intravenous administration of 2,6 dimethoxyphenol or 2,6 dimethoxyquinol in the cat produced a response similar to the delayed toxic effects of M&B 16.573 but not preceded by anaesthesia. The toxic effects of these compounds were

restricted ((arylalkyl)piperidinyl)benzimidazolone analog and study of its activity as ORL1 (NOP) and .kappa.-opioid agonist and its poten
Opioids...

prepn. of (quinolizidinyl)benzimidazolones as conformationally restricted ((arylalkyl)piperidinyl)benzimidazolone analogs and study of their activity as ORL1 (NOP) opioid receptor ligands

Drug discovery...

prepn. of (quinolizidinyl)benzimidazolones as conformationally restricted ((arylalkyl)piperidinyl)benzimidazolone analogs, study of their activity as ORL1 (NOP) opioid receptor ligands and their struc

CAS REGISTRY NUMBERS:

16148-06-4DP 256640-48-9DP conformationally restricted analogs and derivs., prepn. of (quinolizidinyl)benzimidazolones as conformationally restricted ((arylalkyl)piperidinyl)benzimidazolone analogs and study of their activity as ORL1 (NOP) opioid receptor ligands

1501-05-9 5470-11-1 22418-80-0 1493-27-2 prepn. of (quinolizidinyl)benzimidazolones as conformationally restricted ((arylalkyl)piperidinyl)benzimidazolone analogs and study of their activity as ORL1 (NOP) opioid receptor ligands

663609-38-9P SR-14135; prepn. of (quinolizidinyl)benzimidazolones as conformationally restricted ((arylalkyl)piperidinyl)benzimidazolone analogs and study of their activity as ORL1 (NOP) opioid receptor ligands

663609-41-4P SR-14136; prepn. of (quinolizidinyl)benzimidazolone SR-14136 as conformationally restricted ((arylalkyl)piperidinyl)benzimidazolone analog, study of its activity as ORL1 (NOP) and .kappa.-opioid agonist and potential use as analgesic

663609-44-7P SR-14137; prepn. of (quinolizidinyl)benzimidazolones as conformationally restricted ((arylalkyl)piperidinyl)benzimidazolone analogs and study of their activity as ORL1 (NOP) opioid receptor ligands

663609-47-0P SR-14138; prepn. of (quinolizidinyl)benzimidazolones as conformationally restricted ((arylalkyl)piperidinyl)benzimidazolone analogs and study of their activity as ORL1 (NOP) opioid receptor ligands

663609-50-5P SR-14139; prepn. of (quinolizidinyl)benzimidazolones as conformationally restricted ((arylalkyl)piperidinyl)benzimidazolone analogs and study of their activity as ORL1 (NOP) opioid receptor ligands

663609-52-7P SR-14140; prepn. of (quinolizidinyl)benzimidazolones as conformationally restricted ((arylalkyl)piperidinyl)benzimidazolone analogs and study of their activity as ORL1 (NOP) opioid receptor ligands

? ds

Set	Items	Description
S1	457763	MORPHINE OR ATROPINE OR OPIOID OR QUININE (W)ELECTRON (W) - TRANSFER (W) AGENT
S2	7978	QUINOL OR TOCOPHEROL(W)PHOSPHATE
S3	17	S1 AND S2
S4	12	RD (unique items)
S5	12	Sort S4/ALL/PY

? save temp morphine

Temp SearchSave "MORPHINE" stored

? b chemlit

>>> 32 is unauthorized

>>>1 of the specified files is not available

31oct06 08:12:17 User291725 Session D6.3

\$3.92 0.350 DialUnits File73

\$15.50 5 Type(s) in Format 9